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## Chapter 8

# 3D Shape Registration Using Spectral Graph Embedding and Probabilistic Matching

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## Abstract

In this book chapter we address the problem of 3D shape registration and we propose a novel technique based on spectral graph theory and probabilistic matching. Recent advancement in shape acquisition technology has led to the capture of large amounts of 3D data. Existing real-time multi-camera 3D acquisition methods provide a frame-wise reliable visual-hull or mesh representations for real 3D animation sequences. The task of 3D shape analysis involves tracking, recognition, registration, etc. Analyzing 3D data in a single framework is still a challenging task considering the large variability of the data gathered with different acquisition devices. 3D shape registration is one such challenging shape analysis task. The main contribution of this chapter is to extend the spectral graph matching methods to very large graphs by combining spectral graph matching with Laplacian embedding. Since the embedded representation of a graph is obtained by dimensionality reduction we claim that the existing spectral-based methods are not easily applicable. We discuss solutions for the exact and inexact graph isomorphism problems and recall the main spectral properties of the combinatorial graph Laplacian; We provide a novel analysis of the commute-time embedding that allows us to interpret the latter in terms of the PCA of a graph, and to select the appropriate dimension of the associated embedded metric space; We derive a unit hyper-sphere normalization for the commute-time embedding that allows us to register two shapes with different samplings; We propose a novel method to find the eigenvalue-eigenvector ordering and the eigenvector sign using the eigensignature (histogram) which is invariant to the isometric shape deformations and fits well in the spectral graph matching framework, and we present a probabilistic shape matching formulation using an expectation maximization point registration algorithm which alternates between aligning the eigenbases and finding a vertex-to-vertex assignment.

## 8.1 Introduction

In this chapter we discuss the problem of 3D shape registration. Recent advancement in shape acquisition technology has led to the capture of large amounts of 3D data. Existing real-time multi-camera 3D acquisition methods provide a frame-wise reliable visual-hull or mesh representations for real 3D animation sequences [1, 2, 3, 4, 5, 6]. The task of 3D shape analysis involves tracking, recognition, registration, etc. Analyzing 3D data in a single framework is still a challenging task considering the large variability of the data gathered with different acquisition devices. 3D shape registration is one such challenging shape analysis task. The major difficulties in shape registration arise due to: 1) variation in the shape acquisition techniques, 2) local deformations in non-rigid shapes, 3) large acquisition discrepancies (*e.g.*, holes, topology change, surface acquisition noise), 4) local scale change.

Most of the previous attempts of shape matching can be broadly categorized as *extrinsic* or *intrinsic*

approaches depending on how they analyze the properties of the underlying manifold. Extrinsic approaches mainly focus on finding a global or local rigid transformation between two 3D shapes.

There is large set of approaches based on variations of Iterative Closest Point (ICP) algorithm [?, ?, 7] that falls in the category of extrinsic approaches. However, the majority of these approaches compute rigid transformations for shape registration and are not directly applicable to non-rigid shapes. Intrinsic approaches are a natural choice for finding dense correspondences between articulated shapes, as they embed the shape in some canonical domain which preserves some important properties of the manifold, *e.g.*, geodesics and angles. Intrinsic approaches are preferable over extrinsic as they provide a global representation which is invariant to non-rigid deformations that are common in the real-world 3D shapes.

Interestingly, mesh representation also enables the adaptation of well established graph matching algorithms that use eigenvalues and eigenvectors of graph matrices, and are theoretically well investigated in the framework of *Spectral Graph Theory* (SGT) *e.g.*, [8, 9]. Existing methods in SGT are mainly theoretical results applied to small graphs and under the premise that eigenvalues can be computed exactly. However, spectral graph matching does not easily generalize to very large graphs due to the following reasons: 1) eigenvalues are approximately computed using eigen-solvers, 2) eigenvalue multiplicity and hence ordering change are not well studied, 3) exact matching is intractable for very large graphs. It is important to note that these methods mainly focus on exact graph matching while majority of the real-world graph matching applications involve graphs with different cardinality and for which only a subgraph isomorphism can be sought.

The main contribution of this work is to extend the spectral graph methods to very large graphs by combining spectral graph matching with *Laplacian Embedding*. Since the embedded representation of a graph is obtained by dimensionality reduction we claim that the existing SGT methods (*e.g.*, [8]) are not easily applicable. The major contributions of this work are the following: 1) we discuss solutions for the exact and inexact graph isomorphism problems and recall the main spectral properties of the combinatorial graph Laplacian, 2) we provide a novel analysis of the commute-time embedding that allows us to interpret the latter in terms of the PCA of a graph, and to select the appropriate dimension of the associated embedded metric space, 3) we derive a unit hyper-sphere normalization for the commute-time embedding that allows us to register two shapes with different samplings, 4) we propose a novel method to find the eigenvalue-eigenvector ordering and the eigenvector signs using the eigensignatures (histograms) that are invariant to the isometric shape deformations and which fits well in the spectral graph matching framework, 5) we present a probabilistic shape matching formulation using *expectation maximization* for point registration algorithm which alternates between aligning the eigenbases and finding a vertex-to-vertex assignment.

The existing graph matching methods that use intrinsic representations are: [10, 11, 12, 13, 14, 15,

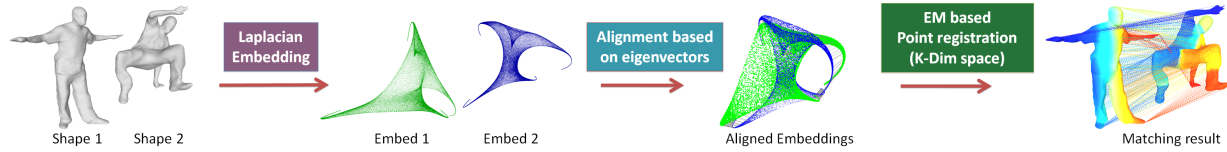


Figure 8.1: Overview of the proposed method. First, a Laplacian embedding is obtained for each shape. Next, these embeddings are aligned to handle the issue of sign flip and ordering change using the histogram matching. Finally, an Expectation-Maximization based point registration is performed to obtain dense probabilistic matching between two shapes.

16, 17]. There is another class of methods that allows to combine intrinsic (geodesics) and extrinsic (appearance) features and which were previously successfully applied for matching features in pairs of images [?, ?, ?, 19, 20, 21, 22, 23, 24, 25]. Some recent approaches apply hierarchical matching to find dense correspondences [26, 27, 28]. However, many of these graph matching algorithms suffer from the problem of either computational intractability or a lack of proper distance metric (*w.r.t.* underlying manifold structure) as the Euclidean metric is not directly applicable while computing distances on non-rigid shapes. A recent benchmarking of shape matching methods was performed in [29]. Recently, a few methods proposed a diffusion framework for the task of shape registration [30, 31, 32].

In this chapter we present an intrinsic approach for unsupervised 3D shape registration first proposed in [14, 33]. In the first step, dimensionality reduction is performed using the graph Laplacian which allows us to embed a 3D shape in an isometric subspace invariant to non-rigid deformations. This leads to an embedded point cloud representation where each vertex of the underlying graph is mapped to a point in a  $K$ -dimensional metric space. Thus, the problem of non-rigid 3D shape registration is transformed into a  $K$ -dimensional point registration task. However, before point registration, the two eigen spaces need to be correctly aligned. This alignment is critical for the spectral matching methods because the two eigen spaces are defined up to the signs and the ordering of the eigenvectors of their Laplacian matrices. This is achieved by a novel matching method that uses histograms of eigenvectors as eigensignatures. In the final step, a point registration method based on a variant of the expectation-maximization (EM) algorithm [34] is applied in order to register two sets of points associated with the Laplacian embeddings of the two shapes. The proposed algorithm alternates between the estimation of an orthogonal transformation matrix associated with the alignment of the two eigen spaces and the computation of probabilistic vertex-to-vertex assignment. Figure 1.1 presents the overview of the proposed method. According to the results summarized in [29], this method is one among the best performing unsupervised shape matching algorithms.

**Chapter Overview:** Graph matrices are introduced in section 1.2. The problem of exact graph isomorphism and existing solutions are discussed in section 1.3. Section 1.4 deals with dimensionality reduction using the graph Laplacian in order to obtain embedded representations for 3D shapes. In the same section we discuss the PCA of graph embeddings and propose a unit hyper-sphere normalization for these embeddings along with a method to choose the embedding dimension. Section 1.5 introduces the formulation of maximum subgraph isomorphism before presenting a two-step method for 3D shape registration. In the first step Laplacian embeddings are aligned using histogram matching while in the second step we briefly discuss an EM point registration method to obtain probabilistic shape registration. Finally we present shape matching results in section 1.6 and conclude with a brief discussion in section 1.7.

## 8.2 Graph Matrices

A shape can be treated as a connected *undirected weighted graph*  $\mathcal{G} = \{\mathcal{V}, \mathcal{E}\}$  where  $\mathcal{V}(\mathcal{G}) = \{v_1, \dots, v_n\}$  is the vertex set,  $\mathcal{E}(\mathcal{G}) = \{e_{ij}\}$  is the edge set. Let  $\mathbf{W}$  be the weighted adjacency matrix of this graph. Each  $(i, j)^{\text{th}}$  entry of  $\mathbf{W}$  matrix stores weight  $w_{ij}$  whenever there is an edge  $e_{ij} \in \mathcal{E}(\mathcal{G})$  between graph vertices  $v_i$  and  $v_j$  and 0 otherwise with all the diagonal elements set to 0. We use the following notations: The degree  $d_i$  of a graph vertex  $d_i = \sum_{i \sim j} w_{ij}$  ( $i \sim j$  denotes the set of vertices  $v_j$  which are adjacent to  $v_i$ ), the *degree matrix*  $\mathbf{D} = \text{diag}[d_1 \dots d_i \dots d_n]$ , the  $n \times 1$  vector  $\mathbb{1} = (1 \dots 1)^{\top}$  (the constant vector), the  $n \times 1$  *degree vector*  $\mathbf{d} = \mathbf{D}\mathbb{1}$ , and the *graph volume*  $\text{Vol}(\mathcal{G}) = \sum_i d_i$ .

In spectral graph theory, it is common [35, 36] to use the following expression for the edge weights:

$$w_{ij} = e^{-\frac{\text{dist}^2(v_i, v_j)}{\sigma^2}}, \quad (8.1)$$

where  $\text{dist}(v_i, v_j)$  denotes any distance metric between two vertices and  $\sigma$  is a free parameter. In the case of a *fully connected graph*, matrix  $\mathbf{W}$  is also referred to as the *similarity matrix*. The *normalized weighted adjacency matrix* writes  $\tilde{\mathbf{W}} = \mathbf{D}^{-1/2} \mathbf{W} \mathbf{D}^{-1/2}$ . The *transition matrix* of the non-symmetric reversible Markov chain associated with the graph is  $\tilde{\mathbf{W}}_R = \mathbf{D}^{-1} \mathbf{W} = \mathbf{D}^{-1/2} \tilde{\mathbf{W}} \mathbf{D}^{1/2}$ .

### 8.2.1 Variants of the Graph Laplacian Matrix

We can now build the concept of the *graph Laplacian operator*. We consider the following variants of the Laplacian matrix [37, 36, 38]:

- The *unnormalized Laplacian* which is also referred to as the *combinatorial Laplacian*  $\mathbf{L}$ ,
- the *normalized Laplacian*  $\tilde{\mathbf{L}}$ , and

- the *random-walk Laplacian*  $\tilde{\mathbf{L}}_R$  also referred to as the *discrete Laplace operator*.

In more detail we have:

$$\mathbf{L} = \mathbf{D} - \mathbf{W} \quad (8.2)$$

$$\tilde{\mathbf{L}} = \mathbf{D}^{-1/2} \mathbf{L} \mathbf{D}^{-1/2} = \mathbf{I} - \tilde{\mathbf{W}} \quad (8.3)$$

$$\tilde{\mathbf{L}}_R = \mathbf{D}^{-1} \mathbf{L} = \mathbf{I} - \tilde{\mathbf{W}}_R \quad (8.4)$$

with the following relations between these matrices:

$$\mathbf{L} = \mathbf{D}^{1/2} \tilde{\mathbf{L}} \mathbf{D}^{1/2} = \mathbf{D} \tilde{\mathbf{L}}_R \quad (8.5)$$

$$\tilde{\mathbf{L}} = \mathbf{D}^{-1/2} \mathbf{L} \mathbf{D}^{-1/2} = \mathbf{D}^{1/2} \tilde{\mathbf{L}}_R \mathbf{D}^{-1/2} \quad (8.6)$$

$$\tilde{\mathbf{L}}_R = \mathbf{D}^{-1/2} \tilde{\mathbf{L}} \mathbf{D}^{1/2} = \mathbf{D}^{-1} \mathbf{L}. \quad (8.7)$$

### 8.3 Spectral Graph Isomorphism

Let  $\mathcal{G}_A$  and  $\mathcal{G}_B$  be two *undirected weighted graphs* with the same number of nodes,  $n$ , and let  $\mathbf{W}_A$  and  $\mathbf{W}_B$  be their adjacency matrices. They are real-symmetric matrices. In the general case, the number  $r$  of distinct eigenvalues of these matrices is smaller than  $n$ . The standard spectral methods only apply to those graphs whose adjacency matrices have  $n$  distinct eigenvalues (each eigenvalue has multiplicity one), which implies that the eigenvalues can be ordered.

Graph isomorphism [39] can be written as the following minimization problem:

$$\mathbf{P}^* = \arg \min_{\mathbf{P}} \|\mathbf{W}_A - \mathbf{P} \mathbf{W}_B \mathbf{P}^\top\|_F^2 \quad (8.8)$$

where  $\mathbf{P}$  is an  $n \times n$  permutation matrix (see appendix A.1) with  $\mathbf{P}^*$  as the desired vertex-to-vertex permutation matrix and  $\|\bullet\|_F$  is the Frobenius norm defined by (see appendix A.2):

$$\|\mathbf{W}\|_F^2 = \langle \mathbf{W}, \mathbf{W} \rangle = \sum_{i=1}^n \sum_{j=1}^n w_{ij}^2 = \text{tr}(\mathbf{W}^\top \mathbf{W}) \quad (8.9)$$

Let:

$$\mathbf{W}_A = \mathbf{U}_A \mathbf{\Lambda}_A \mathbf{U}_A^\top \quad (8.10)$$

$$\mathbf{W}_B = \mathbf{U}_B \mathbf{\Lambda}_B \mathbf{U}_B^\top \quad (8.11)$$

be the eigen-decompositions of the two matrices with  $n$  eigenvalues  $\mathbf{\Lambda}_A = \text{diag}[\alpha_i]$  and  $\mathbf{\Lambda}_B = \text{diag}[\beta_i]$  and  $n$  orthonormal eigenvectors, the column vectors of  $\mathbf{U}_A$  and  $\mathbf{U}_B$ .

### 8.3.1 An Exact Spectral Solution

If there exists a vertex-to-vertex correspondence that makes (1.8) equal to 0, we have:

$$\mathbf{W}_A = \mathbf{P}^* \mathbf{W}_B \mathbf{P}^{*\top}. \quad (8.12)$$

This implies that the adjacency matrices of the two graphs should have the same eigenvalues. Moreover, if the eigenvalues are non null and, the matrices  $\mathbf{U}_A$  and  $\mathbf{U}_B$  have full rank and are uniquely defined by their  $n$  orthonormal column vectors (which are the eigenvectors of  $\mathbf{W}_A$  and  $\mathbf{W}_B$ ), then  $\alpha_i = \beta_i, \forall i, 1 \leq i \leq n$  and  $\mathbf{\Lambda}_A = \mathbf{\Lambda}_B$ . From (1.12) and using the eigen-decompositions of the two graph matrices we obtain:

$$\mathbf{\Lambda}_A = \mathbf{U}_A^\top \mathbf{P}^* \tilde{\mathbf{U}}_B \mathbf{\Lambda}_B \tilde{\mathbf{U}}_B^\top \mathbf{P}^{*\top} \mathbf{U}_A = \mathbf{\Lambda}_B, \quad (8.13)$$

where the matrix  $\tilde{\mathbf{U}}_B$  is defined by:

$$\tilde{\mathbf{U}}_B = \mathbf{U}_B \mathbf{S}. \quad (8.14)$$

Matrix  $\mathbf{S} = \text{diag}[s_i]$ , with  $s_i = \pm 1$ , is referred to as a sign matrix with the property  $\mathbf{S}^2 = \mathbf{I}$ . Post multiplication of  $\mathbf{U}_B$  with a sign matrix takes into account the fact that the eigenvectors (the column vectors of  $\mathbf{U}_B$ ) are only defined up to a sign. Finally we obtain the following permutation matrix:

$$\mathbf{P}^* = \mathbf{U}_B \mathbf{S} \mathbf{U}_A^\top. \quad (8.15)$$

Therefore, one may notice that there are as many solutions as the cardinality of the set of matrices  $\mathbf{S}_n$ , i.e.,  $|\mathbf{S}_n| = 2^n$ , and that *not all of these solutions correspond to a permutation matrix*. This means that there exist some matrices  $\mathbf{S}^*$  that exactly make  $\mathbf{P}^*$  a permutation matrix. Hence, all those permutation matrices that satisfy (1.15) are solutions of the exact graph isomorphism problem. Notice that once the permutation has been estimated, one can write that the rows of  $\mathbf{U}_B$  can be aligned with the rows of  $\mathbf{U}_A$ :

$$\mathbf{U}_A = \mathbf{P}^* \mathbf{U}_B \mathbf{S}^*. \quad (8.16)$$

The rows of  $\mathbf{U}_A$  and of  $\mathbf{U}_B$  can be interpreted as isometric embeddings of the two graph vertices: A vertex  $v_i$  of  $\mathcal{G}_A$  has as coordinates the  $i^{\text{th}}$  row of  $\mathbf{U}_A$ . This means that the spectral graph isomorphism problem becomes a point registration problem, where graph vertices are represented by points in  $\mathbb{R}^n$ . To conclude, the exact graph isomorphism problem has a spectral solution based on: (i) the eigen-decomposition of the two graph matrices, (ii) the ordering of their eigenvalues, and (iii) the choice of a sign for each eigenvector.

### 8.3.2 The Hoffman-Wielandt Theorem

The Hoffman-Wielandt theorem [40, 41] is the fundamental building block of spectral graph isomorphism. The theorem holds for normal matrices; Here, we restrict the analysis to real symmetric matrices, although the generalization to Hermitian matrices is straightforward:



**Theorem 1**

(Hoffman and Wielandt) If  $\mathbf{W}_A$  and  $\mathbf{W}_B$  are real-symmetric matrices, and if  $\alpha_i$  and  $\beta_i$  are their eigenvalues arranged in increasing order,  $\alpha_1 \leq \dots \leq \alpha_i \leq \dots \leq \alpha_n$  and  $\beta_1 \leq \dots \leq \beta_i \leq \dots \leq \beta_n$ , then

$$\sum_{i=1}^n (\alpha_i - \beta_i)^2 \leq \|\mathbf{W}_A - \mathbf{W}_B\|_F^2. \quad (8.17)$$

*Proof:* The proof is derived from [9, 42]. Consider the eigen-decompositions of matrices  $\mathbf{W}_A$  and  $\mathbf{W}_B$ , (1.10), (1.11). Notice that for the time being we are free to prescribe the ordering of the eigenvalues  $\alpha_i$  and  $\beta_i$  and hence the ordering of the column vectors of matrices  $\mathbf{U}_A$  and  $\mathbf{U}_B$ . By combining (1.10) and (1.11) we write:

$$\mathbf{U}_A \mathbf{\Lambda}_A \mathbf{U}_A^\top - \mathbf{U}_B \mathbf{\Lambda}_B \mathbf{U}_B^\top = \mathbf{W}_A - \mathbf{W}_B \quad (8.18)$$

or, equivalently:

$$\mathbf{\Lambda}_A \mathbf{U}_A^\top \mathbf{U}_B - \mathbf{U}_A^\top \mathbf{U}_B \mathbf{\Lambda}_B = \mathbf{U}_A^\top (\mathbf{W}_A - \mathbf{W}_B) \mathbf{U}_B. \quad (8.19)$$

By the unitary-invariance of the Frobenius norm (see appendix A.2 ) and with the notation  $\mathbf{Z} = \mathbf{U}_A^\top \mathbf{U}_B$  we obtain:

$$\|\mathbf{\Lambda}_A \mathbf{Z} - \mathbf{Z} \mathbf{\Lambda}_B\|_F^2 = \|\mathbf{W}_A - \mathbf{W}_B\|_F^2, \quad (8.20)$$

which is equivalent to:

$$\sum_{i=1}^n \sum_{j=1}^n (\alpha_i - \beta_j)^2 z_{ij}^2 = \|\mathbf{W}_A - \mathbf{W}_B\|_F^2. \quad (8.21)$$

The coefficients  $x_{ij} = z_{ij}^2$  can be viewed as the entries of a doubly-stochastic matrix  $\mathbf{X}$ :  $x_{ij} \geq 0, \sum_{i=1}^n x_{ij} = 1, \sum_{j=1}^n x_{ij} = 1$ . Using these properties, we obtain:

$$\begin{aligned} \sum_{i=1}^n \sum_{j=1}^n (\alpha_i - \beta_j)^2 z_{ij}^2 &= \sum_{i=1}^n \alpha_i^2 + \sum_{j=1}^n \beta_j^2 - 2 \sum_{i=1}^n \sum_{j=1}^n z_{ij}^2 \alpha_i \beta_j \\ &\geq \sum_{i=1}^n \alpha_i^2 + \sum_{j=1}^n \beta_j^2 - 2 \max_Z \left\{ \sum_{i=1}^n \sum_{j=1}^n z_{ij}^2 \alpha_i \beta_j \right\}. \end{aligned} \quad (8.22)$$

Hence, the minimization of (1.21) is equivalent to the maximization of the last term in (1.22). We can modify our maximization problem to admit all the doubly-stochastic matrices. In this way we seek an extremum over a convex compact set. The maximum over this compact set is larger than or equal to our maximum:

$$\max_{Z \in O_n} \left\{ \sum_{i=1}^n \sum_{j=1}^n z_{ij}^2 \alpha_i \beta_j \right\} \leq \max_{X \in \mathcal{D}_n} \left\{ \sum_{i=1}^n \sum_{j=1}^n x_{ij} \alpha_i \beta_j \right\} \quad (8.23)$$

where  $O_n$  is the set of orthogonal matrices and  $\mathcal{D}_n$  is the set of doubly stochastic matrices (see appendix A.1). Let  $c_{ij} = \alpha_i \beta_j$  and hence one can write that the right term in the equation above as the dot-product of two matrices:

$$\langle \mathbf{X}, \mathbf{C} \rangle = \text{tr}(\mathbf{X}\mathbf{C}) = \sum_{i=1}^n \sum_{j=1}^n x_{ij} c_{ij}. \quad (8.24)$$

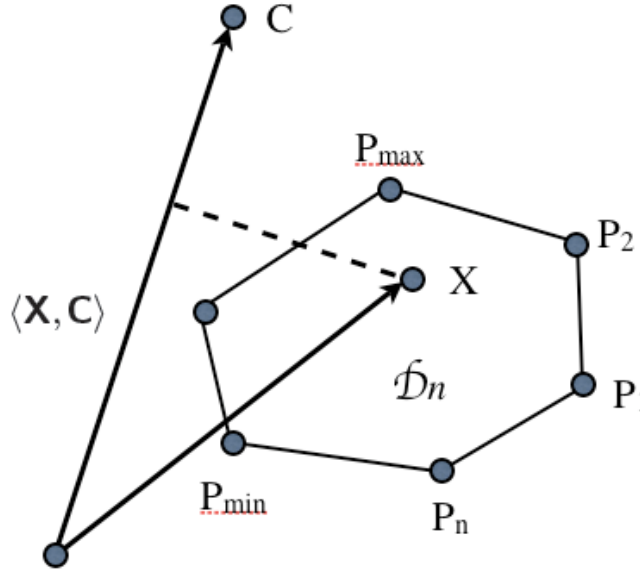


Figure 8.2: This figure illustrates the maximization of the dot-product  $\langle \mathbf{X}, \mathbf{C} \rangle$ . The two matrices can be viewed as vectors of dimension  $n^2$ . Matrix  $\mathbf{X}$  belongs to a compact convex set whose extreme points are the permutation matrices  $\mathbf{P}_1, \mathbf{P}_2, \dots, \mathbf{P}_n$ . Therefore, the projection of this set (i.e.,  $\mathcal{D}_n$ ) onto  $\mathbf{C}$  has projected permutation matrices at its extremes, namely  $\langle \mathbf{P}_{\min}, \mathbf{X} \rangle$  and  $\langle \mathbf{P}_{\max}, \mathbf{X} \rangle$  in this example.

Therefore, this expression can be interpreted as the projection of  $\mathbf{X}$  onto  $\mathbf{C}$ , see figure 1.2. The Birkhoff theorem (appendix A.1) tells us that the set  $\mathcal{D}_n$  of doubly stochastic matrices is a compact convex set. We obtain that the extrema (minimum and maximum) of the projection of  $\mathbf{X}$  onto  $\mathbf{C}$  occur at the projections of one of the extreme points of this convex set, which correspond to permutation matrices. Hence, the maximum of  $\langle \mathbf{X}, \mathbf{C} \rangle$  is  $\langle \mathbf{P}_{\max}, \mathbf{X} \rangle$  and we obtain:

$$\max_{X \in \mathcal{D}_n} \left\{ \sum_{i=1}^n \sum_{j=1}^n x_{ij} \alpha_i \beta_j \right\} = \sum_{i=1}^n \alpha_i \beta_{\pi(i)}. \quad (8.25)$$

By substitution in (1.22) we obtain:

$$\sum_{i=1}^n \sum_{j=1}^n (\alpha_i - \beta_j)^2 z_{ij}^2 \geq \sum_{i=1}^n (\alpha_i - \beta_{\pi(i)})^2. \quad (8.26)$$

If the eigenvalues are in increasing order then the permutation that satisfies theorem 1.17 is the identity matrix, i.e.,  $\pi(i) = i$ . Indeed, let's assume that for some indices  $k$  and  $k+1$  we have:  $\pi(k) = k+1$  and  $\pi(k+1) = k$ . Since  $\alpha_k \leq \alpha_{k+1}$  and  $\beta_k \leq \beta_{k+1}$ , the following inequality holds:

$$(\alpha_k - \beta_k)^2 + (\alpha_{k+1} - \beta_{k+1})^2 \leq (\alpha_k - \beta_{k+1})^2 + (\alpha_{k+1} - \beta_k)^2 \quad (8.27)$$

and hence (1.17) holds. ■

**Corollary 1.1**

The inequality (1.17) becomes an equality when the eigenvectors of  $\mathbf{W}_A$  are aligned with the eigenvectors of  $\mathbf{W}_B$  up to a sign ambiguity:

$$\mathbf{U}_B = \mathbf{U}_A \mathbf{S}. \quad (8.28)$$

*Proof:* Since the minimum of (1.21) is achieved for  $\mathbf{X} = \mathbf{I}$  and since the entries of  $\mathbf{X}$  are  $z_{ij}^2$ , we have that  $z_{ii} = \pm 1$ , which corresponds to  $\mathbf{Z} = \mathbf{S}$ . ■

**Corollary 1.2**

If  $\mathbf{Q}$  is an orthogonal matrix, then

$$\sum_{i=1}^n (\alpha_i - \beta_i)^2 \leq \|\mathbf{W}_A - \mathbf{QW}_B\mathbf{Q}^\top\|_F^2. \quad (8.29)$$

*Proof:* Since the eigen-decomposition of matrix  $\mathbf{QW}_B\mathbf{Q}^\top$  is  $(\mathbf{QU}_B)\mathbf{\Lambda}_B(\mathbf{QU}_B)^\top$  and since it has the same eigenvalues as  $\mathbf{W}_B$ , the inequality (1.29) holds and hence corollary 1.2. ■

These corollaries will be useful in the case of spectral graph matching methods presented below.

**8.3.3 Umeyama's Method**

The exact spectral matching solution presented in section 1.3.1 finds a permutation matrix satisfying (1.15). This requires an exhaustive search over the space of all possible  $2^n$  matrices. Umeyama's method presented in [8] proposes a relaxed solution to this problem as outlined below.

Umeyama [8] addresses the problem of *weighted graph matching* within the framework of spectral graph theory. He proposes two methods, the first for *undirected weighted graphs* and the second for *directed weighted graphs*. The adjacency matrix is used in both cases. Let's consider the case of undirected graphs. The eigenvalues are (possibly with multiplicities):

$$\mathbf{W}_A: \alpha_1 \leq \dots \leq \alpha_i \leq \dots \leq \alpha_n \quad (8.30)$$

$$\mathbf{W}_B: \beta_1 \leq \dots \leq \beta_i \leq \dots \leq \beta_n. \quad (8.31)$$

**Theorem 2**

(Umeyama) If  $\mathbf{W}_A$  and  $\mathbf{W}_B$  are real-symmetric matrices with  $n$  distinct eigenvalues (that can be ordered),  $\alpha_1 < \dots < \alpha_i < \dots < \alpha_n$  and  $\beta_1 < \dots < \beta_i < \dots < \beta_n$ , the minimum of :

$$J(\mathbf{Q}) = \|\mathbf{W}_A - \mathbf{QW}_B\mathbf{Q}^\top\|_F^2 \quad (8.32)$$

is achieved for:

$$\mathbf{Q}^* = \mathbf{U}_A \mathbf{S} \mathbf{U}_B^\top \quad (8.33)$$

and hence (1.29) becomes an equality:

$$\sum_{i=1}^n (\alpha_i - \beta_i)^2 = \|\mathbf{W}_A - \mathbf{Q}^* \mathbf{W}_B \mathbf{Q}^{*\top}\|_F^2. \quad (8.34)$$

*Proof:* The proof is straightforward. By corollary 1.2, the Hoffman-Wielandt theorem applies to matrices  $\mathbf{W}_A$  and  $\mathbf{Q} \mathbf{W}_B \mathbf{Q}^\top$ . By corollary 1.1, the equality (1.34) is achieved for:

$$\mathbf{Z} = \mathbf{U}_A^\top \mathbf{Q}^* \mathbf{U}_B = \mathbf{S} \quad (8.35)$$

and hence (1.33) holds. ■

Notice that (1.33) can be written as:

$$\mathbf{U}_A = \mathbf{Q}^* \mathbf{U}_B \mathbf{S} \quad (8.36)$$

which is a *relaxed* version of (1.16): The permutation matrix in the exact isomorphism case is replaced by an orthogonal matrix.

**A Heuristic for Spectral Graph Matching:** Let us consider again the exact solution outlined in section 1.3.1. Umeyama suggests a heuristic in order to avoid exhaustive search over all possible  $2^n$  matrices that satisfy (1.15). One may easily notice that:

$$\|\mathbf{P} - \mathbf{U}_A \mathbf{S} \mathbf{U}_B^\top\|_F^2 = 2n - 2\text{tr}(\mathbf{U}_A \mathbf{S} (\mathbf{P} \mathbf{U}_B)^\top). \quad (8.37)$$

Using Umeyama's notations,  $\bar{\mathbf{U}}_A = [|u_{ij}|]$ ,  $\bar{\mathbf{U}}_B = [|v_{ij}|]$  (the entries of  $\bar{\mathbf{U}}_A$  are the absolute values of the entries of  $\mathbf{U}_A$ ), one may further notice that:

$$\text{tr}(\mathbf{U}_A \mathbf{S} (\mathbf{P} \mathbf{U}_B)^\top) = \sum_{i=1}^n \sum_{j=1}^n s_j u_{ij} v_{\pi(i)j} \leq \sum_{i=1}^n \sum_{j=1}^n |u_{ij}| |v_{\pi(i)j}| = \text{tr}(\bar{\mathbf{U}}_A \bar{\mathbf{U}}_B^\top \mathbf{P}^\top). \quad (8.38)$$

The minimization of (1.37) is equivalent to the maximization of (1.38) and the maximal value that can be attained by the latter is  $n$ . Using the fact that both  $\mathbf{U}_A$  and  $\mathbf{U}_B$  are orthogonal matrices, one can easily conclude that:

$$\text{tr}(\bar{\mathbf{U}}_A \bar{\mathbf{U}}_B^\top \mathbf{P}^\top) \leq n. \quad (8.39)$$

Umeyama concludes that when the two graphs are isomorphic, the optimum permutation matrix maximizes  $\text{tr}(\bar{\mathbf{U}}_A \bar{\mathbf{U}}_B^\top \mathbf{P}^\top)$  and this can be solved by the Hungarian algorithm [43].

When the two graphs are not exactly isomorphic, theorem 1 and theorem 2 allow us to relax the permutation matrices to the group of orthogonal matrices. Therefore with similar arguments as above we obtain:

$$\text{tr}(\mathbf{U}_A \mathbf{S} \mathbf{U}_B^\top \mathbf{Q}^\top) \leq \text{tr}(\bar{\mathbf{U}}_A \bar{\mathbf{U}}_B^\top \mathbf{Q}^\top) \leq n. \quad (8.40)$$

The permutation matrix obtained with the Hungarian algorithm can be used as an initial solution that can then be improved by some hill-climbing or relaxation technique [8].

The spectral matching solution presented in this section is not directly applicable to large graphs. In the next section we introduce the notion of dimensionality reduction for graphs which will lead to a tractable graph matching solution.

## 8.4 Graph Embedding and Dimensionality Reduction

For *large* and *sparse* graphs, the results of section 1.3 and Umeyama's method (section 1.3.3) hold only *weakly*. Indeed, one cannot guarantee that all the eigenvalues have multiplicity equal to one: the presence of symmetries causes some of eigenvalues to have an algebraic multiplicity greater than one. Under these circumstances and due to numerical approximations, it might not be possible to properly order the eigenvalues. Moreover, for very large graphs with thousands of vertices it is not practical to compute all its eigenvalue-eigenvector pairs. This means that one has to devise a method that is able to match shapes using a small set of eigenvalues and eigenvectors.

One elegant way to overcome this problem, is to reduce the dimension of the eigenspace, along the line of spectral dimensionality reductions techniques. The eigendecomposition of graph Laplacian matrices (introduced in section 1.2.1) is a popular choice for the purpose of dimensionality reduction [35].

### 8.4.1 Spectral Properties of the Graph Laplacian

The spectral properties of the Laplacian matrices introduced in section 1.2.1 have been thoroughly studied. They are summarized in table 1.1. We derive some subtle properties of the combinatorial Laplacian which

Laplacian	Null space	Eigenvalues	Eigenvectors
$\mathbf{L} = \mathbf{U}\mathbf{\Lambda}\mathbf{U}^\top$	$\mathbf{u}_1 = \mathbb{1}$	$0 = \lambda_1 < \lambda_2 \leq \dots \leq \lambda_n$	$\mathbf{u}_{i>1}^\top \mathbb{1} = 0, \mathbf{u}_i^\top \mathbf{u}_j = \delta_{ij}$
$\tilde{\mathbf{L}} = \tilde{\mathbf{U}}\mathbf{\Gamma}\tilde{\mathbf{U}}^\top$	$\tilde{\mathbf{u}}_1 = \mathbf{D}^{1/2}\mathbb{1}$	$0 = \gamma_1 < \gamma_2 \leq \dots \leq \gamma_n$	$\tilde{\mathbf{u}}_{i>1}^\top \mathbf{D}^{1/2}\mathbb{1} = 0, \tilde{\mathbf{u}}_i^\top \tilde{\mathbf{u}}_j = \delta_{ij}$
$\tilde{\mathbf{L}}_R = \mathbf{T}\mathbf{\Gamma}\mathbf{T}^{-1}, \mathbf{T} = \mathbf{D}^{-1/2}\tilde{\mathbf{U}}$	$\mathbf{t}_1 = \mathbb{1}$	$0 = \gamma_1 < \gamma_2 \leq \dots \leq \gamma_n$	$\mathbf{t}_{i>1}^\top \mathbf{D}\mathbb{1} = 0, \mathbf{t}_i^\top \mathbf{D}\mathbf{t}_j = \delta_{ij}$

Table 8.1: Summary of the spectral properties of the Laplacian matrices. Assuming a connected graph, the null eigenvalue  $(\lambda_1, \gamma_1)$  has multiplicity one. The first non null eigenvalue  $(\lambda_2, \gamma_2)$  is known as the Fiedler value and its multiplicity is, in general, equal to one. The associated eigenvector is denoted the Fiedler vector [37].

will be useful for the task of shape registration. In particular, we will show that the eigenvectors of the

combinatorial Laplacian can be interpreted as directions of maximum variance (principal components) of the associated embedded shape representation. We note that the embeddings of the normalized and random-walk Laplacians have different spectral properties which make them less interesting for shape registration, i.e., Appendix A.3.

**The combinatorial Laplacian.** Let  $\mathbf{L} = \mathbf{U}\mathbf{\Lambda}\mathbf{U}^\top$  be the spectral decomposition of the combinatorial Laplacian with  $\mathbf{U}\mathbf{U}^\top = \mathbf{I}$ . Let  $\mathbf{U}$  be written as:

$$\mathbf{U} = \begin{bmatrix} u_{11} & \dots & u_{1k} & \dots & u_{1n} \\ \vdots & & \vdots & & \vdots \\ u_{n1} & \dots & u_{nk} & \dots & u_{nn} \end{bmatrix} \quad (8.41)$$

Each column of  $\mathbf{U}$ ,  $\mathbf{u}_k = (u_{1k} \dots u_{ik} \dots u_{nk})^\top$  is an eigenvector associated with the eigenvalue  $\lambda_k$ . From the definition of  $\mathbf{L}$  in (1.2) (see [35]) one can easily see that  $\lambda_1 = 0$  and that  $\mathbf{u}_1 = \mathbb{1}$  (a constant vector). Hence,  $\mathbf{u}_{k \geq 2}^\top \mathbb{1} = 0$  and by combining this with  $\mathbf{u}_k^\top \mathbf{u}_k = 1$ , we derive the following proposition:

**Proposition 1**

*The components of the non-constant eigenvectors of the combinatorial Laplacian satisfy the following constraints:*

$$\sum_{i=1}^n u_{ik} = 0, \quad \forall k, 2 \leq k \leq n \quad (8.42)$$

$$-1 < u_{ik} < 1, \quad \forall i, k, 1 \leq i \leq n, 2 \leq k \leq n. \quad (8.43)$$

Assuming a connected graph,  $\lambda_1$  has multiplicity equal to one [36]. Let's organize the eigenvalues of  $\mathbf{L}$  in increasing order:  $0 = \lambda_1 < \lambda_2 \leq \dots \leq \lambda_n$ . We prove the following proposition [37]:

**Proposition 2**

*For all  $k \leq n$ , we have  $\lambda_k \leq 2 \max_i(d_i)$ , where  $d_i$  is the degree of vertex  $i$ .*

*Proof:* The largest eigenvalue of  $\mathbf{L}$  corresponds to the maximization of the Rayleigh quotient, or

$$\lambda_n = \max_{\mathbf{u}} \frac{\mathbf{u}^\top \mathbf{L} \mathbf{u}}{\mathbf{u}^\top \mathbf{u}}. \quad (8.44)$$

We have  $\mathbf{u}^\top \mathbf{L} \mathbf{u} = \sum_{e_{ij}} w_{ij} (u_i - u_j)^2$ . From the inequality  $(a - b)^2 \leq 2(a^2 + b^2)$  we obtain:

$$\lambda_n \leq \frac{2 \sum_{e_{ij}} w_{ij} (u_i^2 + u_j^2)}{\sum_i u_i^2} = \frac{2 \sum_i d_i u_i^2}{\sum_i u_i^2} \leq 2 \max_i(d_i). \quad \blacksquare \quad (8.45)$$

This ensures an upper limit on the eigenvalues of  $\mathbf{L}$ . By omitting the zero eigenvalue and associated eigenvector, we can rewrite  $\mathbf{L}$  as:

$$\mathbf{L} = \sum_{k=2}^n \lambda_k \mathbf{u}_k \mathbf{u}_k^\top. \quad (8.46)$$

Each entry  $u_{ik}$  of an eigenvector  $\mathbf{u}_k$  can be interpreted as a real-valued function that projects a graph vertex  $v_i$  onto that vector. The mean and variance of the set  $\{u_{ik}\}_{i=1}^n$  are therefore a measure of how the graph *spreads* when projected onto the  $k$ -th eigenvector. This is clarified by the following result:

**Proposition 3**

*The mean  $\bar{u}_k$  and the variance  $\sigma_{u_k}$  of an eigenvector  $\mathbf{u}_k$ . For  $2 \leq k \leq n$ , and  $1 \leq i \leq n$  we have*

$$\bar{u}_k = \sum_{i=1}^n u_{ik} = 0 \quad (8.47)$$

$$\sigma_{u_k} = \frac{1}{n} \sum_{i=1}^n (u_{ik} - \bar{u}_k)^2 = \frac{1}{n} \quad (8.48)$$

*Proof:* These results can be easily obtained from  $\mathbf{u}_{k \geq 2}^\top \mathbf{1} = 0$  and  $\mathbf{u}_k^\top \mathbf{u}_k = 1$ . ■

These properties will be useful while aligning two Laplacian embeddings and thus registering two 3D shapes.

### 8.4.2 Principal Component Analysis of a Graph Embedding

The Moore-Penrose pseudo-inverse of the Laplacian can be written as:

$$\begin{aligned} \mathbf{L}^\dagger &= \mathbf{U} \mathbf{\Lambda}^{-1} \mathbf{U}^\top \\ &= (\mathbf{\Lambda}^{-\frac{1}{2}} \mathbf{U}^\top)^\top (\mathbf{\Lambda}^{-\frac{1}{2}} \mathbf{U}^\top) \\ &= \mathbf{X}^\top \mathbf{X} \end{aligned} \quad (8.49)$$

where  $\mathbf{\Lambda}^{-1} = \text{diag}(0, 1/\lambda_2, \dots, 1/\lambda_n)$ .

The symmetric semi-definite positive matrix  $\mathbf{L}^\dagger$  is a *Gram* matrix with the same eigenvectors as those of the graph Laplacian. When omitting the null eigenvalue and associated constant eigenvector,  $\mathbf{X}$  becomes a  $(n-1) \times n$  matrix whose columns are the coordinates of the graph's vertices in an *embedded (or feature) space*, i.e.,  $\mathbf{X} = [\mathbf{x}_1 \dots \mathbf{x}_j \dots \mathbf{x}_n]$ . It is interesting to note that the entries of  $\mathbf{L}^\dagger$  may be viewed as *kernel* dot-products, or a Gram matrix [44]. The Gram-matrix representation allows us to embed the graph in an Euclidean feature-space where each vertex  $v_j$  of the graph is a feature point represented as  $\mathbf{x}_j$ .

The left pseudo-inverse operator of the Laplacian  $\mathbf{L}$ , satisfying  $\mathbf{L}^\dagger \mathbf{L} \mathbf{u} = \mathbf{u}$  for any  $\mathbf{u} \perp \text{null}(\mathbf{L})$ , is also called the *Green function* of the heat equation. Under the assumption that the graph is connected and thus  $\mathbf{L}$

has an eigenvalue  $\lambda_1 = 0$  with multiplicity 1, we obtain:

$$\mathbf{L}^\dagger = \sum_{k=2}^n \frac{1}{\lambda_k} \mathbf{u}_k \mathbf{u}_k^\top. \quad (8.50)$$

The Green function is intimately related to random walks on graphs, and can be interpreted probabilistically as follows. Given a Markov chain such that each graph vertex is the state, and the transition from vertex  $v_i$  is possible to any adjacent vertex  $v_j \sim v_i$  with probability  $w_{ij}/d_i$ , the expected number of steps required to reach vertex  $v_j$  from  $v_i$ , called the *access* or *hitting time*  $O(v_i, v_j)$ . The expected number of steps in a round trip from  $v_i$  to  $v_j$  is called the *commute-time distance*:  $\text{CTD}^2(v_i, v_j) = O(v_i, v_j) + O(v_j, v_i)$ . The commute-time distance [45] can be expressed in terms of the entries of  $\mathbf{L}^\dagger$ :

$$\begin{aligned} \text{CTD}^2(v_i, v_j) &= \text{Vol}(\mathcal{G})(\mathbf{L}^\dagger(i, i) + \mathbf{L}^\dagger(j, j) - 2\mathbf{L}^\dagger(i, j)) \\ &= \text{Vol}(\mathcal{G}) \left( \sum_{k=2}^n \frac{1}{\lambda_k} \mathbf{u}_{ik}^2 + \sum_{k=2}^n \frac{1}{\lambda_k} \mathbf{u}_{jk}^2 - 2 \sum_{k=2}^n \frac{1}{\lambda_k} \mathbf{u}_{ik} \mathbf{u}_{jk} \right) \\ &= \text{Vol}(\mathcal{G}) \sum_{k=2}^n \left( \lambda_k^{-1/2} (\mathbf{u}_{ik} - \mathbf{u}_{jk}) \right)^2 \\ &= \text{Vol}(\mathcal{G}) \|\mathbf{x}_i - \mathbf{x}_j\|^2, \end{aligned} \quad (8.51)$$

where the volume of the graph,  $\text{Vol}(\mathcal{G})$  is the sum of the degrees of all the graph vertices. The CTD function is positive-definite and sub-additive, thus defining a *metric* between the graph vertices, referred to as *commute-time* (or *resistance*) *distance* [46]. The CTD is inversely related to the number and length of paths connecting two vertices. Unlike the shortest-path (geodesic) distance, CTD captures the connectivity structure of the graph volume rather than a single path between the two vertices. The great advantage of the commute-time distance over the shortest geodesic path is that it is robust to topological changes and therefore is well suited for characterizing complex shapes. Since the volume is a graph constant, we obtain:

$$\text{CTD}^2(v_i, v_j) \propto \|\mathbf{x}_i - \mathbf{x}_j\|^2. \quad (8.52)$$

Hence, the Euclidean distance between any two feature points  $\mathbf{x}_i$  and  $\mathbf{x}_j$  is the commute time distance between the graph vertex  $v_i$  and  $v_j$ .

Using the first  $K$  non-null eigenvalue-eigenvector pairs of the Laplacian  $\mathbf{L}$ , the *commute-time embedding* of the graph's nodes corresponds to the column vectors of the  $K \times n$  matrix  $\mathbf{X}$ :

$$\mathbf{X}_{K \times n} = \mathbf{\Lambda}_K^{-1/2} (\mathbf{U}_{n \times K})^\top = [\mathbf{x}_1 \dots \mathbf{x}_j \dots \mathbf{x}_n]. \quad (8.53)$$

From (1.43) and (1.53) one can easily infer lower and upper bounds for the  $i$ -th coordinate of  $\mathbf{x}_j$ :

$$-\lambda_i^{-1/2} < x_{ji} < \lambda_i^{-1/2}. \quad (8.54)$$



The last equation implies that the graph embedding stretches along the eigenvectors with a factor that is inversely proportional to the square root of the eigenvalues. Theorem 3 below characterizes the smallest non-null  $K$  eigenvalue-eigenvector pairs of  $\mathbf{L}$  as the directions of maximum variance (the principal components) of the commute-time embedding.

### Theorem 3

*The largest eigenvalue-eigenvector pairs of the pseudo-inverse of the combinatorial Laplacian matrix are the principal components of the commute-time embedding, i.e., the points  $\mathbf{X}$  are zero-centered and have a diagonal covariance matrix.*

*Proof:* Indeed, from (1.47) we obtain a zero-mean while from (1.53) we obtain a diagonal covariance matrix:

$$\bar{\mathbf{x}} = \frac{1}{n} \sum_{i=1}^n \mathbf{x}_i = \frac{1}{n} \mathbf{\Lambda}^{-\frac{1}{2}} \begin{pmatrix} \sum_{i=1}^n \mathbf{u}_{i2} \\ \vdots \\ \sum_{i=1}^n \mathbf{u}_{ik+1} \end{pmatrix} = \begin{pmatrix} 0 \\ \vdots \\ 0 \end{pmatrix} \quad (8.55)$$

$$\Sigma_X = \frac{1}{n} \mathbf{X} \mathbf{X}^\top = \frac{1}{n} \mathbf{\Lambda}^{-\frac{1}{2}} \mathbf{U}^\top \mathbf{U} \mathbf{\Lambda}^{-\frac{1}{2}} = \frac{1}{n} \mathbf{\Lambda}^{-1} \quad (8.56)$$

■.

Figure 1.3 shows the projection of graph (in this case 3D shape represented as meshes) vertices on eigenvectors.

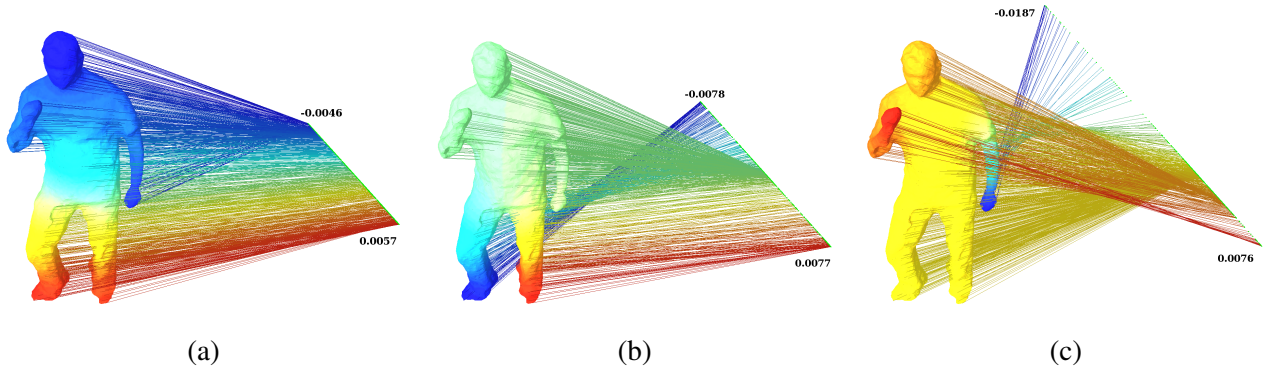


Figure 8.3: This is an illustration of the concept of the PCA of a graph embedding. The graph's vertices are projected onto the second, third and fourth eigenvectors of the Laplacian matrix. These eigenvectors can be viewed as the principal directions of the shape.

### 8.4.3 Choosing the Dimension of the Embedding

A direct consequence of theorem 3 is that the embedded graph representation is centered and the eigenvectors of the combinatorial Laplacian are the directions of maximum variance. The *principal* eigenvectors correspond to the eigenvectors associated with the  $K$  largest eigenvalues of the  $\mathbf{L}^\dagger$ , i.e.,  $\lambda_2^{-1} \geq \lambda_3^{-1} \geq \dots \geq \lambda_K^{-1}$ . The variance along vector  $\mathbf{u}_k$  is  $\lambda_k^{-1}/n$ . Therefore, the total variance can be computed from the trace of the  $\mathbf{L}^\dagger$  matrix :

$$\text{tr}(\mathbf{\Sigma}_X) = \frac{1}{n} \text{tr}(\mathbf{L}^\dagger). \quad (8.57)$$

A standard way of choosing the principal components is to use the *scree diagram*:

$$\theta(K) = \frac{\sum_{k=2}^{K+1} \lambda_k^{-1}}{\sum_{k=2}^n \lambda_k^{-1}}. \quad (8.58)$$

The selection of the first  $K$  principal eigenvectors therefore depends on the spectral fall-off of the inverses of the eigenvalues. In spectral graph theory, the dimension  $K$  is chosen on the basis of the existence of an eigengap, such that  $\lambda_{K+2} - \lambda_{K+1} > t$  with  $t > 0$ . In practice it is extremely difficult to find such an eigengap, in particular in the case of sparse graphs that correspond to a discretized manifold. Instead, we propose to select the dimension of the embedding in the following way. Notice that (1.58) can be written as  $\theta(K) = A/(A+B)$  with  $A = \sum_{k=2}^{K+1} \lambda_k^{-1}$  and  $B = \sum_{k=K+2}^n \lambda_k^{-1}$ . Moreover, from the fact that the  $\lambda_k$ 's are arranged in increasing order, we obtain  $B \leq (n-K-1)\lambda_{K+1}^{-1}$ . Hence:

$$\theta_{\min} \leq \theta(K) \leq 1, \quad (8.59)$$

with

$$\theta_{\min} = \frac{\sum_{k=2}^{K+1} \lambda_k^{-1}}{\sum_{k=2}^K \lambda_k^{-1} + (n-K)\lambda_{K+1}^{-1}}. \quad (8.60)$$

This lower bound can be computed from the  $K$  smallest non null eigenvalues of the combinatorial Laplacian matrix. Hence, one can choose  $K$  such that the sum of the first  $K$  eigenvalues of the  $\mathbf{L}^\dagger$  matrix is a good approximation of the total variance, e.g.,  $\theta_{\min} = 0.95$ .

### 8.4.4 Unit Hyper-sphere Normalization

One disadvantage of the standard embeddings is that, when two shapes have large difference in sampling the embeddings will differ by a significant scale factor. In order to avoid this we can *normalize* the embedding such that the vertex coordinates lie on a unit sphere of dimension  $K$ , which yields:

$$\hat{\mathbf{x}}_i = \frac{\mathbf{x}_i}{\|\mathbf{x}_i\|}. \quad (8.61)$$

In more detail, the  $k$ -th coordinate of  $\hat{\mathbf{x}}_i$  writes as:

$$\hat{\mathbf{x}}_{ik} = \frac{\lambda_k^{-\frac{1}{2}} \mathbf{u}_{ik}}{\left( \sum_{l=2}^{K+1} \lambda_l^{-\frac{1}{2}} \mathbf{u}_{il}^2 \right)^{1/2}}. \quad (8.62)$$

## 8.5 Spectral Shape Matching

In the previous sections we discussed solutions for the exact and inexact graph isomorphism problems, we recalled the main spectral properties of the combinatorial graph Laplacian, and we provided a novel analysis of the commute-time embedding that allows to interpret the latter in terms of the PCA of a graph, and to select the appropriate dimension  $K \ll n$  of the associated embedded metric space. In this section we address the problem of 3D shape registration and we illustrate how the material developed above can be exploited in order to build a robust algorithm for spectral shape matching.

Let's consider two shapes described by two graphs,  $\mathcal{G}_A$  and  $\mathcal{G}_B$  where  $|\mathcal{V}_A| = n$  and  $|\mathcal{V}_B| = m$ . Let  $\mathbf{L}_A$  and  $\mathbf{L}_B$  be their corresponding graph Laplacians. Without loss of generality, one can choose the same dimension  $K \ll \min(n, m)$  for the two embeddings. This yields the following eigen decompositions:

$$\mathbf{L}_A = \mathbf{U}_{n \times K} \mathbf{\Lambda}_K (\mathbf{U}_{n \times K})^\top \quad (8.63)$$

$$\mathbf{L}_B = \mathbf{U}'_{m \times K} \mathbf{\Lambda}'_K (\mathbf{U}'_{m \times K})^\top. \quad (8.64)$$

For each one of these graphs, one can build two *isomorphic* embedded representations, as follows:

- An *unnormalized Laplacian embedding* that uses the  $K$  rows of  $\mathbf{U}_{n \times K}$  as the Euclidean coordinates of the vertices of  $\mathcal{G}_A$  (as well as the  $K$  rows of  $\mathbf{U}'_{m \times K}$  as the Euclidean coordinates of the vertices of  $\mathcal{G}_B$ ), and
- A *normalized commute-time embedding* given by (1.61), i.e.,  $\hat{\mathbf{X}}_A = [\hat{\mathbf{x}}_1 \dots \hat{\mathbf{x}}_j \dots \hat{\mathbf{x}}_n]$  (as well as  $\hat{\mathbf{X}}_B = [\hat{\mathbf{x}}'_1 \dots \hat{\mathbf{x}}'_j \dots \hat{\mathbf{x}}'_m]$ ). We recall that each column  $\hat{\mathbf{x}}_j$  (and respectively  $\hat{\mathbf{x}}'_j$ ) is a  $K$ -dimensional vector corresponding to a vertex  $v_j$  of  $\mathcal{G}_A$  (and respectively  $v'_j$  of  $\mathcal{G}_B$ ).

### 8.5.1 Maximum Subgraph Matching and Point Registration

Let's apply the graph isomorphism framework of Section 1.3 to the two graphs. They are embedded into two congruent spaces of dimension  $\mathbb{R}^K$ . If the smallest  $K$  non-null eigenvalues associated with the two embeddings are distinct and can be ordered, i.e.:

$$\lambda_2 < \dots < \lambda_k < \dots < \lambda_{K+1} \quad (8.65)$$

$$\lambda'_2 < \dots < \lambda'_k < \dots < \lambda'_{K+1} \quad (8.66)$$

then, the Umeyama method could be applied. If one uses the unnormalized Laplacian embeddings just defined, (1.33) becomes:

$$\mathbf{Q}^* = \mathbf{U}_{n \times K} \mathbf{S}_K (\mathbf{U}'_{m \times K})^\top \quad (8.67)$$

Notice that here the sign matrix  $\mathbf{S}$  defined in 1.33 became a  $K \times K$  matrix denoted by  $\mathbf{S}_K$ . We now assume that the eigenvalues  $\{\lambda_2, \dots, \lambda_{K+1}\}$  and  $\{\lambda'_2, \dots, \lambda'_{K+1}\}$  *cannot be reliably ordered*. This can be modeled by multiplication with a  $K \times K$  permutation matrix  $\mathbf{P}_K$ :

$$\mathbf{Q} = \mathbf{U}_{n \times K} \mathbf{S}_K \mathbf{P}_K (\mathbf{U}'_{m \times K})^\top \quad (8.68)$$

Pre-multiplication of  $(\mathbf{U}'_{m \times K})^\top$  with  $\mathbf{P}_K$  permutes its rows such that  $\mathbf{u}'_k \rightarrow \mathbf{u}'_{\pi(k)}$ . Each entry  $q_{ij}$  of the  $n \times m$  matrix  $\mathbf{Q}$  can therefore be written as:

$$q_{ij} = \sum_{k=2}^{K+1} s_k u_{ik} u'_{j\pi(k)} \quad (8.69)$$

Since both  $\mathbf{U}_{n \times K}$  and  $\mathbf{U}'_{m \times K}$  are column-orthonormal matrices, the dot-product defined by (1.69) is equivalent to the cosine of the angle between two  $K$ -dimensional vectors. This means that each entry of  $\mathbf{Q}$  is such that  $-1 \leq q_{ij} \leq +1$  and that two vertices  $v_i$  and  $v'_j$  are matched if  $q_{ij}$  is close to 1.

One can also use the normalized commute-time coordinates and define an equivalent expression as above:

$$\hat{\mathbf{Q}} = \hat{\mathbf{X}}^\top \mathbf{S}_K \mathbf{P}_K \hat{\mathbf{X}}' \quad (8.70)$$

with:

$$\hat{q}_{ij} = \sum_{k=2}^{K+1} s_k \hat{x}_{ik} \hat{x}'_{j\pi(k)} \quad (8.71)$$

Because both sets of points  $\hat{\mathbf{X}}$  and  $\hat{\mathbf{X}}'$  lie on a  $K$ -dimensional unit hyper-sphere, we also have  $-1 \leq \hat{q}_{ij} \leq +1$ .

It should however be emphasized that the rank of the  $n \times m$  matrices  $\mathbf{Q}, \hat{\mathbf{Q}}$  is equal to  $K$ . Therefore, these matrices cannot be viewed as *relaxed permutation matrices* between the two graphs. In fact they define many-to-many correspondences between the vertices of the first graph and the vertices of the second graph, this being due to the fact that the graphs are embedded on a low-dimensional space. This is one of the main differences between our method proposed in the next section and the Umeyama method, as well as many other subsequent methods, that use all eigenvectors of the graph. As it will be explained below, our formulation leads to a shape matching method that will alternate between aligning their eigenbases and finding a vertex-to-vertex assignment.

It is possible to extract a one-to-one assignment matrix from  $\mathbf{Q}$  (or from  $\hat{\mathbf{Q}}$ ) using either dynamic programming or an assignment method technique such as the Hungarian algorithm. Notice that this assignment is conditioned by the choice of a sign matrix  $\mathbf{S}_K$  and of a permutation matrix  $\mathbf{P}_K$ , i.e.,  $2^K K!$  possibilities,

and that not all these choices correspond to a valid sub-isomorphism between the two graphs. Let's consider the case of the normalized commute-time embedding; there is an equivalent formulation for the unnormalized Laplacian embedding. The two graphs are described by two sets of points,  $\hat{\mathbf{X}}$  and  $\hat{\mathbf{X}}'$ , both lying onto the  $K$ -dimensional unity hyper-sphere. The  $K \times K$  matrix  $\mathbf{S}_K \mathbf{P}_K$  transforms one graph embedding onto the other graph embedding. Hence, one can write  $\hat{\mathbf{x}}_i = \mathbf{S}_K \mathbf{P}_K \hat{\mathbf{x}}'_j$  if vertex  $v_i$  matches  $v_j$ . More generally Let  $\mathbf{R}_K = \mathbf{S}_K \mathbf{P}_K$  and let's extend the domain of  $\mathbf{R}_K$  to all possible orthogonal matrices of size  $K \times K$ , namely  $\mathbf{R}_K \in O_K$  or the *orthogonal group* of dimension  $K$ . We can now write the following criterion whose minimization over  $\mathbf{R}_K$  guarantees an optimal solution for registering the vertices of the first graph with the vertices of the second graph:

$$\min_{\mathbf{R}_K} \sum_{i=1}^n \sum_{j=1}^m \hat{q}_{ij} \|\hat{\mathbf{x}}_i - \mathbf{R}_K \hat{\mathbf{x}}'_j\|^2 \quad (8.72)$$

One way to solve minimization problems such as (1.72) is to use a point registration algorithm that alternates between (i) estimating the  $K \times K$  orthogonal transformation  $\mathbf{R}_K$ , which aligns the  $K$ -dimensional coordinates associated with the two embeddings, and (ii) updating the assignment variables  $\hat{q}_{ij}$ . This can be done using either ICP-like methods (the  $\hat{q}_{ij}$ 's are binary variables), or EM-like methods (the  $\hat{q}_{ij}$ 's are posterior probabilities of assignment variables). As we just outlined above, matrix  $\mathbf{R}_K$  belongs to the orthogonal group  $O_K$ . Therefore this framework differs from standard implementations of ICP and EM algorithms that usually estimate a 2-D or 3-D *rotation* matrix which belong to the *special orthogonal group*.

It is well established that ICP algorithms are easily trapped in local minima. The EM algorithm recently proposed in [34] is able to converge to a good solution starting with a rough initial guess and is robust to the presence of outliers. Nevertheless, the algorithm proposed in [34] performs well under *rigid transformations* (rotation and translation), whereas in our case we have to estimate a more general orthogonal transformation that incorporates both rotations and reflections. Therefore, before describing in detail an EM algorithm well suited for solving the problem at hand, we discuss the issue of estimating an initialization for the transformation aligning the  $K$  eigenvectors of the first embedding with those of the second embedding and we propose a practical method for initializing this transformation (namely, matrices  $\mathbf{S}_K$  and  $\mathbf{P}_K$  in (1.70)) based on comparing the histograms of these eigenvectors, or *eigensignatures*.

### 8.5.2 Aligning Two Embeddings Based on Eigensignatures

Both the unnormalized Laplacian embedding and the normalized commute-time embedding of a graph are represented in a metric space spanned by the eigenvectors of the Laplacian matrix, namely the  $n$ -dimensional vectors  $\{\mathbf{u}_2, \dots, \mathbf{u}_k, \dots, \mathbf{u}_{K+1}\}$ , where  $n$  is the number of graph vertices. They correspond to *eigenfunctions* and each such eigenfunction maps the graph's vertices onto the real line. More precisely, the  $k$ -th

eigenfunction maps a vertex  $v_i$  onto  $u_{ik}$ . Propositions 1 and 3 revealed interesting statistics of the sets  $\{u_{1k}, \dots, u_{ik}, \dots, u_{nk}\}_{k=2}^{K+1}$ . Moreover, theorem 3 provided an interpretation of the eigenvectors in terms of principal directions of the embedded shape. One can therefore conclude that the probability distribution of the components of an eigenvector have interesting properties that make them suitable for comparing two shapes, namely  $-1 < u_{ik} < +1$ ,  $\bar{u}_k = 1/n \sum_{i=1}^n u_{ik} = 0$ , and  $\sigma_k = 1/n \sum_{i=1}^n u_{ik}^2 = 1/n$ . This means that one can build a histogram for each eigenvector and that all these histograms share the same bin width  $w$  and the same number of bins  $b$  [47]:

$$w_k = \frac{3.5\sigma_k}{n^{1/3}} = \frac{3.5}{n^{4/3}} \quad (8.73)$$

$$b_k = \frac{\sup_i u_{ik} - \inf_i u_{ik}}{w_k} \approx \frac{n^{4/3}}{2}. \quad (8.74)$$

We claim that these histograms are eigenvector signatures which are invariant under graph isomorphism. Indeed, let's consider the Laplacian  $\mathbf{L}$  of a shape and we apply the isomorphic transformation  $\mathbf{PLP}^\top$  to this shape, where  $\mathbf{P}$  is a permutation matrix. If  $\mathbf{u}$  is an eigenvector of  $\mathbf{L}$ , it follows that  $\mathbf{Pu}$  is an eigenvector of  $\mathbf{PLP}^\top$  and therefore, while the order of the components of  $\mathbf{u}$  are affected by this transformation, their frequency and hence their probability distribution remain the same. Hence, one may conclude that such a histogram may well be viewed as an *eigensignature*.

We denote with  $H\{\mathbf{u}\}$  the histogram formed with the components of  $\mathbf{u}$  and let  $C(H\{\mathbf{u}\}, H\{\mathbf{u}'\})$  be a similarity measure between two histograms. From the eigenvector properties just outlined, it is straightforward to notice that  $H\{\mathbf{u}\} \neq H\{-\mathbf{u}\}$ : These two histograms are mirror symmetric. Hence, the histogram is not invariant to the sign of an eigenvector. Therefore one can use the eigenvectors' histograms to estimate both the permutation matrix  $\mathbf{P}_K$  and the sign matrix  $\mathbf{S}_K$  in (1.70). The problem of finding one-to-one assignments  $\{\mathbf{u}_k \leftrightarrow s_k \mathbf{u}'_{\pi(k)}\}_{k=2}^{K+1}$  between the two sets of eigenvectors associated with the two shapes is therefore equivalent to the problem of finding one-to-one assignments between their histograms.

Let  $\mathbf{A}_K$  be an assignment matrix between the histograms of the first shape and the histograms of the second shape. Each entry of this matrix is defined by:

$$a_{kl} = \sup[C(H\{\mathbf{u}_k\}, H\{\mathbf{u}'_l\}); C(H\{\mathbf{u}_k\}, H\{-\mathbf{u}'_l\})] \quad (8.75)$$

Similarly, we define a matrix  $\mathbf{B}_K$  that accounts for the *sign assignments*:

$$b_{kl} = \begin{cases} +1 & \text{if } C(H\{\mathbf{u}_k\}, H\{\mathbf{u}'_l\}) \geq C(H\{\mathbf{u}_k\}, H\{-\mathbf{u}'_l\}) \\ -1 & \text{if } C(H\{\mathbf{u}_k\}, H\{\mathbf{u}'_l\}) < C(H\{\mathbf{u}_k\}, H\{-\mathbf{u}'_l\}) \end{cases} \quad (8.76)$$

Extracting a permutation matrix  $\mathbf{P}_K$  from  $\mathbf{A}_K$  is an instance of the bipartite maximum matching problem and the Hungarian algorithm is known to provide an optimal solution to this assignment problem [43].

Moreover, one can use the estimated  $\mathbf{P}_K$  to extract a sign matrix  $\mathbf{S}_K$  from  $\mathbf{B}_K$ . Algorithm 1 estimates an alignment between two embeddings.

---

**Algorithm 1** *Alignment of Two Laplacian Embeddings*

---

**input** : Histograms associated with eigenvectors  $\{\mathbf{u}_k\}_{k=2}^{K+1}$  and  $\{\mathbf{u}'_k\}_{k=2}^{K+1}$ .

**output** : A permutation matrix  $\mathbf{P}_K$  and a sign matrix  $\mathbf{S}_K$ .

- 1: Compute the assignment matrices  $\mathbf{A}_K$  and  $\mathbf{B}_K$ .
  - 2: Compute  $\mathbf{P}_K$  from  $\mathbf{A}_K$  using the Hungarian algorithm.
  - 3: Compute the sign matrix  $\mathbf{S}_K$  using  $\mathbf{P}_K$  and  $\mathbf{B}_K$ .
- 

Figure 1.4 illustrates the utility of the histogram of eigenvectors as eigensignatures for solving the problem of sign flip and change in eigenvector ordering by computing histogram matching. It is interesting to observe that a threshold on the histogram matching score (1.75) allows us to discard the eigenvectors with low similarity cost. Hence, starting with large  $K$  obtained using (1.60), we can limit the number of eigenvectors to just a few, which will be suitable for EM based point registration algorithm proposed in the next section.

### 8.5.3 An EM Algorithm for Shape Matching

As explained in section 1.5.1, the maximum subgraph matching problem reduces to a point registration problem in  $K$  dimensional metric space spanned by the eigenvectors of graph Laplacian where two shapes are represented as point clouds. The initial alignment of Laplacian embeddings can be obtained by matching the histogram of eigenvectors as described in the previous section. In this section we propose an EM algorithm for 3D shape matching that computes a probabilistic vertex-to-vertex assignment between two shapes. The proposed method alternates between the step to estimate an orthogonal transformation matrix associated with the alignment of the two shape embeddings and the step to compute a point-to-point probabilistic assignment variable.

The method is based on a parametric probabilistic model, namely maximum likelihood with missing data. Let us consider the Laplacian embedding of two shapes, i.e., (1.53) :  $\hat{\mathbf{X}} = \{\hat{\mathbf{x}}_i\}_{i=1}^n, \hat{\mathbf{X}}' = \{\hat{\mathbf{x}}'_j\}_{j=1}^m$ , with  $\hat{\mathbf{X}}, \hat{\mathbf{X}}' \subset \mathbb{R}^K$ . Without loss of generality, we assume that the points in the first set,  $\hat{\mathbf{X}}$  are cluster centers of a Gaussian mixture model (GMM) with  $n$  clusters and an additional uniform component that accounts for outliers and unmatched data. The matching  $\hat{\mathbf{X}} \leftrightarrow \hat{\mathbf{X}}'$  will consist in fitting the Gaussian mixture to the set  $\hat{\mathbf{X}}'$ .

Let this Gaussian mixture undergo a  $K \times K$  transformation  $\mathbf{R}$  (for simplicity, we omit the index  $K$ ) with  $\mathbf{R}^\top \mathbf{R} = \mathbf{I}_K, \det(\mathbf{R}) = \pm 1$ , more precisely  $\mathbf{R} \in O_K$ , the group of orthogonal matrices acting on  $\mathbb{R}^K$ . Hence, each cluster in the mixture is parametrized by a prior  $p_i$ , a cluster mean  $\mu_i = \mathbf{R}\hat{\mathbf{x}}_i$ , and a covariance matrix

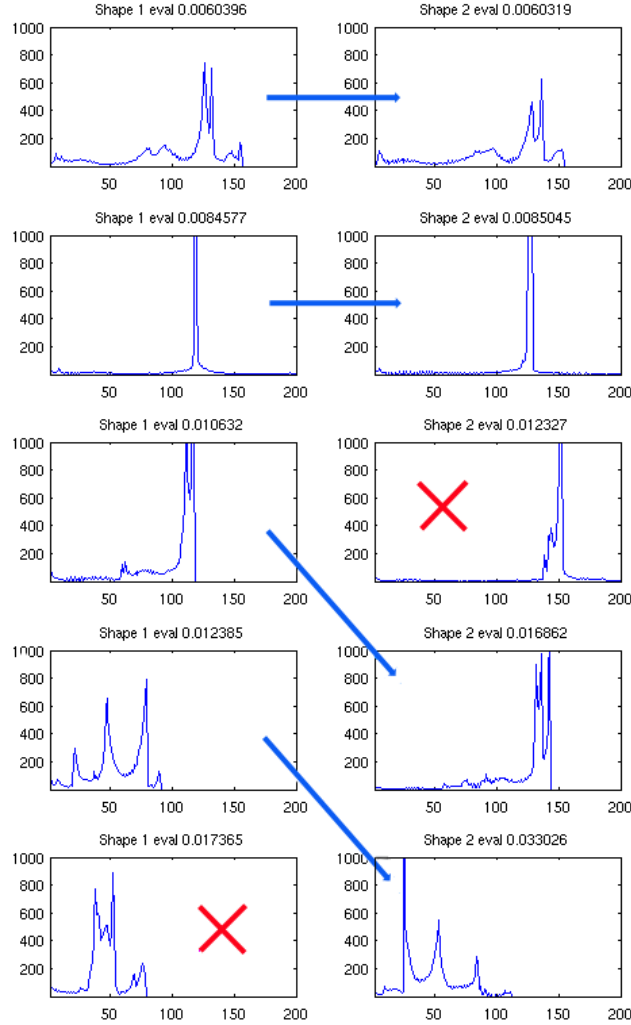


Figure 8.4: An illustration of applicability of eigenvector histogram as eigensignature to detect sign flip and eigenvector ordering change. The blue line shows matched eigenvector pairs and the red-cross depicts discarded eigenvectors.

$\Sigma_i$ . It will be assumed that all the clusters in the mixture have the same priors,  $\{p_i = \pi_{\text{in}}\}_{i=1}^n$ , and the same isotropic covariance matrix,  $\{\Sigma_i = \sigma \mathbf{I}_K\}_{i=1}^n$ . This parametrization leads to the following *observed-data log-likelihood* (with  $\pi_{\text{out}} = 1 - n\pi_{\text{in}}$  and  $\mathcal{U}$  is the uniform distribution):

$$\log P(\mathbf{X}') = \sum_{j=1}^m \log \left( \sum_{i=1}^n \pi_{\text{in}} \mathcal{N}(\mathbf{x}'_j | \mu_i, \sigma) + \pi_{\text{out}} \mathcal{U} \right) \quad (8.77)$$

It is well known that the direct maximization of (1.77) is not tractable and it is more practical to maximize the *expected complete-data log-likelihood* using the EM algorithm, where “complete-data” refers to both the



observed data (the points  $\hat{\mathbf{X}}'$ ) and the missing data (the data-to-cluster assignments). In our case, the above expectation writes (see [34] for details):

$$\mathcal{E}(\mathbf{R}, \sigma) = -\frac{1}{2} \sum_{j=1}^m \sum_{i=1}^n \alpha_{ji} (\|\hat{\mathbf{x}}'_j - \mathbf{R}\hat{\mathbf{x}}_i\|^2 + k \log \sigma), \quad (8.78)$$

where  $\alpha_{ji}$  denotes the posterior probability of an assignment:  $\hat{\mathbf{x}}'_j \leftrightarrow \hat{\mathbf{x}}_i$ :

$$\alpha_{ji} = \frac{\exp(-\|\hat{\mathbf{x}}'_j - \mathbf{R}\hat{\mathbf{x}}_i\|^2 / 2\sigma)}{\sum_{q=1}^n \exp(-\|\hat{\mathbf{x}}'_j - \mathbf{R}\hat{\mathbf{x}}_q\|^2 / 2\sigma) + \emptyset \sigma^{k/2}}, \quad (8.79)$$

where  $\emptyset$  is a constant term associated with the uniform distribution  $\mathcal{U}$ . Notice that one easily obtains the posterior probability of a data point to remain unmatched,  $\alpha_{jn+1} = 1 - \sum_{i=1}^n \alpha_{ij}$ . This leads to the shape matching procedure outlined in Algorithm 2.

---

**Algorithm 2** *EM for shape matching*

---

**input** : Two embedded shapes  $\hat{\mathbf{X}}$  and  $\hat{\mathbf{X}}'$ ;

**output** : Dense correspondences  $\hat{\mathbf{X}} \leftrightarrow \hat{\mathbf{X}}'$  between the two shapes;

- 1: *Initialization*: Set  $\mathbf{R}^{(0)} = \mathbf{S}_K \mathbf{P}_K$  choose a large value for the variance  $\sigma^{(0)}$ ;
- 2: *E-step*: Compute the current posteriors  $\alpha_{ij}^{(q)}$  from the current parameters using (1.79);
- 3: *M-step*: Compute the new transformation  $\mathbf{R}^{(q+1)}$  and the new variance  $\sigma^{(q+1)}$  using the current posteriors:

$$\begin{aligned} \mathbf{R}^{(q+1)} &= \arg \min_{\mathbf{R}} \sum_{i,j} \alpha_{ij}^{(q)} \|\mathbf{x}'_j - \mathbf{R}\mathbf{x}_i\|^2 \\ \sigma^{(q+1)} &= \sum_{i,j} \alpha_{ij}^{(q)} \|\hat{\mathbf{x}}'_j - \mathbf{R}^{(q+1)}\hat{\mathbf{x}}_i\|^2 / k \sum_{i,j} \alpha_{ij}^{(q)} \end{aligned}$$

- 4: *MAP*: Accept the assignment  $\hat{\mathbf{x}}'_j \leftrightarrow \hat{\mathbf{x}}_i$  if  $\max_i \alpha_{ij}^{(q)} > 0.5$ .
- 

## 8.6 Experiments and Results

We have performed several 3D shape registration experiments to evaluate the proposed method. In the first experiment, 3D shape registration is performed on 138 high-resolution (10K-50K vertices) triangular meshes from the publicly available TOSCA dataset [29]. The dataset includes 3 shape classes (human, dog, horse) with simulated transformations. Transformations are split into 9 classes (isometry, topology, small and big holes, global and local scaling, noise, shot noise, sampling). Each transformation class appears in five different strength levels. An estimate of average geodesic distance to ground truth correspondence was computed for performance evaluation (see [29] for details).

We evaluate our method in two settings. In the first setting SM1 we use the commute-time embedding (1.53) while in the second setting SM2 we use the unit hyper-sphere normalized embedding (1.61).

Transform	Strength									
	1		$\leq 2$		$\leq 3$		$\leq 4$		$\leq 5$	
	SM1	SM2	SM1	SM2	SM1	SM2	SM1	SM2	SM1	SM2
<i>Isometry</i>	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
<i>Topology</i>	6.89	5.96	7.92	6.76	7.92	7.14	8.04	7.55	8.41	8.13
<i>Holes</i>	7.32	5.17	8.39	5.55	9.34	6.05	9.47	6.44	12.47	10.32
<i>Micro holes</i>	0.37	0.68	0.39	0.70	0.44	0.79	0.45	0.79	0.49	0.83
<i>Scale</i>	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
<i>Local scale</i>	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
<i>Sampling</i>	11.43	10.51	13.32	12.08	15.70	13.65	18.76	15.58	22.63	19.17
<i>Noise</i>	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
<i>Shot noise</i>	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00	0.00
<b>Average</b>	2.88	2.48	3.34	2.79	3.71	3.07	4.08	3.37	4.89	4.27

Table 8.2: 3D shape registration error estimates (average geodesic distance to ground truth correspondences) using proposed spectral matching method with commute-time embedding (SM1) and unit hyper-sphere normalized embedding (SM2).

Table 1.2 shows the error estimates for dense shape matching using proposed spectral matching method. In the case of some transforms, the proposed method yields zero error because the two meshes had identical triangulations. Figure 1.5 shows some matching results. The colors emphasize the correct matching of body parts while we show only 5% of matches for better visualization. In Figure 1.5(e) the two shapes have large difference in the sampling rate. In this case the matching near the shoulders is not fully correct since we used the commute-time embedding.

Table 1.3 summarizes the comparison of proposed spectral matching method (SM1 and SM2) with generalized multidimensional scaling (GMDS) based matching algorithm introduced in [17] and the Laplace-Beltrami matching algorithm proposed in [10] with two settings LB1 (uses graph Laplacian) and LB2 (uses cotangent weights). GMDS computes correspondence between two shapes by trying to embed one shape into another with minimum distortion. LB1 and LB2 algorithms combines the surface descriptors based on the eigendecomposition of the Laplace-Beltrami operator and the geodesic distances measured on the shapes when calculating the correspondence quality. The above results in a quadratic optimization problem formulation for correspondence detection, and its minimizer is the best possible correspondence. The proposed method clearly outperform the other two methods with minimum average error estimate computed over all the transformations in the dataset.

<b>Method</b>	<b>Strength</b>				
	1	$\leq 2$	$\leq 3$	$\leq 4$	$\leq 5$
<i>LB1</i>	10.61	15.48	19.01	23.22	23.88
<i>LB2</i>	15.51	18.21	22.99	25.26	28.69
<i>GMDS</i>	39.92	36.77	35.24	37.40	39.10
<i>SM1</i>	2.88	3.34	3.71	4.08	4.89
<i>SM2</i>	2.48	2.79	3.07	3.37	4.27

Table 8.3: Average shape registration error estimates over all transforms (average geodesic distance to ground truth correspondences) computed using proposed methods (SM1 and SM2), GMDS [17] and LB1, LB2 [10].

<b>Transform</b>	<b>Strength</b>		
	1	$\leq 3$	$\leq 5$
<i>Isometry</i>	SM1,SM2	SM1,SM2	SM1,SM2
<i>Topology</i>	SM2	SM2	SM2
<i>Holes</i>	SM2	SM2	SM2
<i>Micro holes</i>	SM1	SM1	SM1
<i>Scale</i>	SM1,SM2	SM1,SM2	SM1,SM2
<i>Local scale</i>	SM1,SM2	SM1,SM2	SM1,SM2
<i>Sampling</i>	LB1	SM2	LB2
<i>Noise</i>	SM1,SM2	SM1,SM2	SM1,SM2
<i>Shot noise</i>	SM1,SM2	SM1,SM2	SM1,SM2
<b>Average</b>	SM1,SM2	SM1,SM2	SM1,SM2

Table 8.4: 3D shape registration performance comparison: The proposed methods (SM1 and SM2) performed best by providing minimum average shape registration error over all the transformation classes with different strength as compare to GMDS [17] and LB1, LB2 [10] methods.

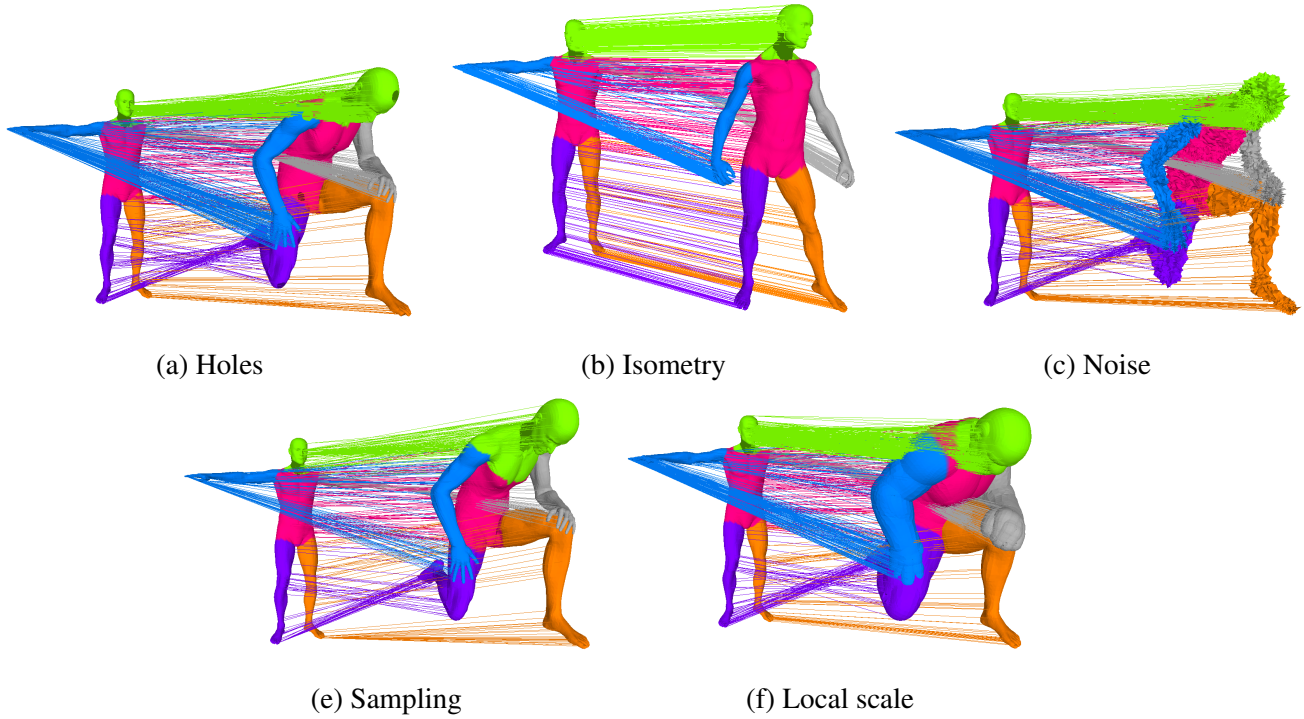


Figure 8.5: 3D shape registration in the presence of different transforms.

In table 1.4, we show a detailed comparison of proposed method with other methods. For a detailed quantitative comparison refer to [29]. The proposed method inherently uses diffusion geometry as opposed to geodesic metric used by other two methods and hence outperform them.

In the second experiment we perform shape registration on two different shapes with similar topology. In Figure 1.6, results of shape registration on different shapes is presented. Figure 1.6(a),(c) shows the initialization step of EM algorithm while Figure 1.6(b),(d) shows the dense matching obtained after EM convergence.

Finally, we show shape matching results on two different human meshes captured with multi-camera system at MIT [5] and University of Surrey [2] in Figure 1.7

## 8.7 Discussion

This chapter describes a 3D shape registration approach that computes dense correspondences between two articulated objects. We address the problem using spectral matching and unsupervised point registration method. We formally introduce graph isomorphism using the Laplacian matrix, and we provide an analysis of the matching problem when the number of nodes in the graph is very large, *i.e.* of the order of  $O(10^4)$ . We show that there is a simple equivalence between graph isomorphism and point registration under the

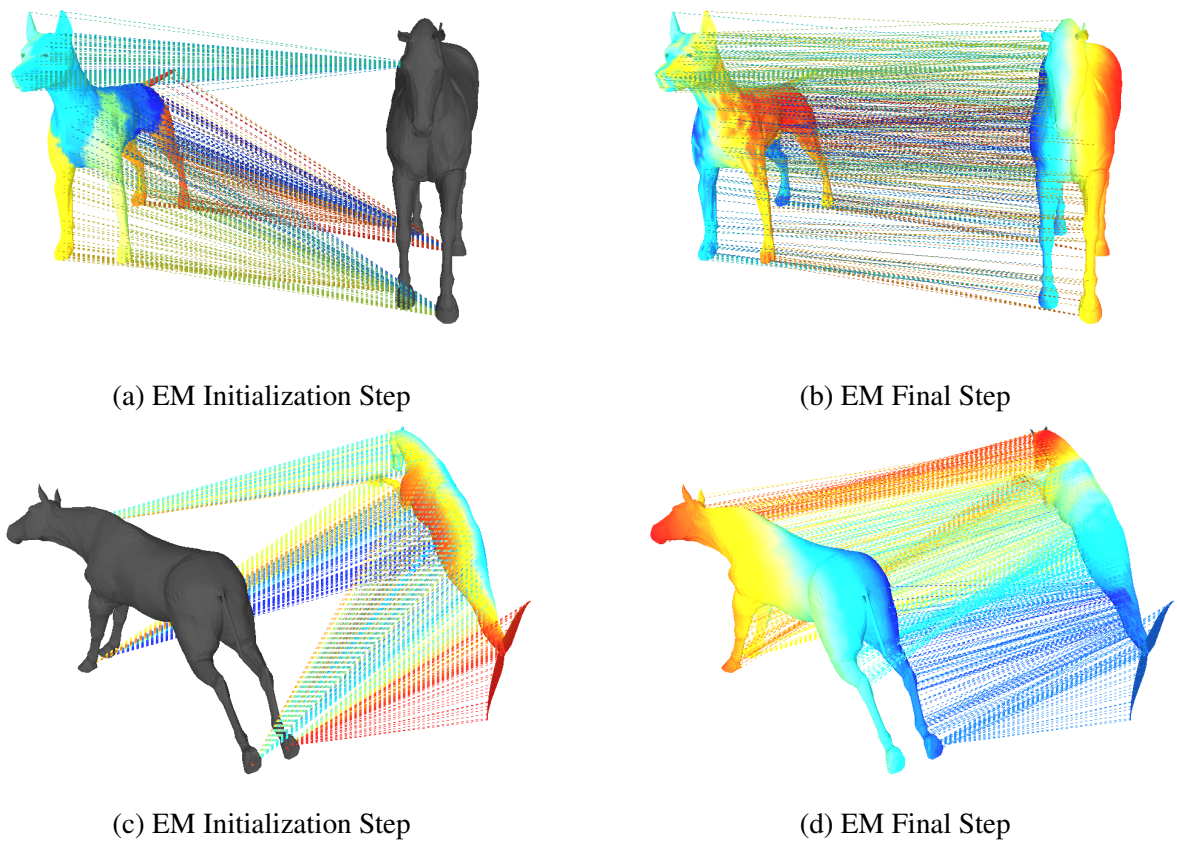


Figure 8.6: 3D shape registration performed on different shapes with similar topology.

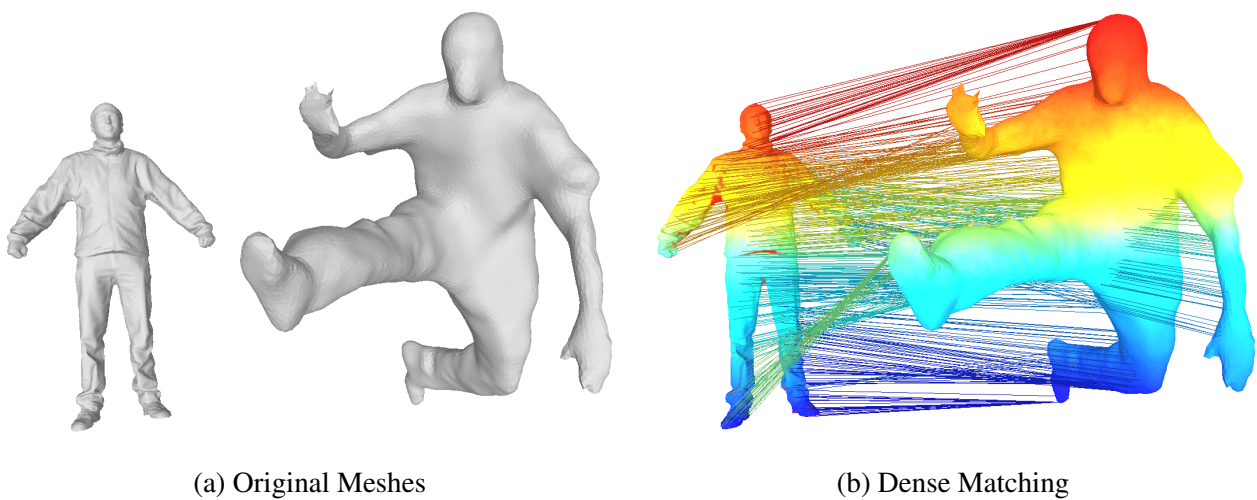


Figure 8.7: 3D shape registration performed on two real meshes captured from different sequence.

group of orthogonal transformations, when the dimension of the embedding space is much smaller than the cardinality of the point-sets.

The eigenvalues of a large sparse Laplacian cannot be reliably ordered. We propose an elegant alternative to eigenvalue ordering, using eigenvector histograms and alignment based on comparing these histograms. The point registration that results from eigenvector alignment yields an excellent initialization for the EM algorithm, subsequently used only to refine the registration.

However, the method is susceptible to large topology changes that might occur in the multi-camera shape acquisition setup due to self-occlusion (originated from complex kinematics poses) and shadow effects. This is because Laplacian embedding is a global representation and any major topology change will lead to large changes in embeddings causing failure of this method. Recently, a new shape registration method proposed in [32] provides robustness to the large topological changes using the heat kernel framework.



# Appendix A

## A.1 Permutation and Doubly-stochastic Matrices

A matrix  $\mathbf{P}$  is called a *permutation* matrix if exactly one entry in each row and column is equal to 1, and all other entries are 0. Left multiplication of a matrix  $\mathbf{A}$  by a permutation matrix  $\mathbf{P}$  permutes the *rows* of  $\mathbf{A}$ , while right multiplication permutes the *columns* of  $\mathbf{A}$ .

Permutation matrices have the following properties:  $\det(\mathbf{P}) = \pm 1$ ,  $\mathbf{P}^\top = \mathbf{P}^{-1}$ , the identity is a permutation matrix, and the product of two permutation matrices is a permutation matrix. Hence the set of permutation matrices  $\mathbf{P} \in \mathcal{P}_n$  constitute a subgroup of the subgroup of orthogonal matrices, denoted by  $O_n$ , and  $\mathcal{P}_n$  has finite cardinality  $n!$ .

A non-negative matrix  $\mathbf{A}$  is a matrix such that all its entries are non-negative. A non-negative matrix with the property that all its row sums are +1 is said to be a (*row*) *stochastic matrix*. A *column stochastic matrix* is the transpose of a row stochastic matrix. A stochastic matrix  $\mathbf{A}$  with the property that  $\mathbf{A}^\top$  is also stochastic is said to be *doubly stochastic*: all row and column sums are +1 and  $a_{ij} \geq 0$ . The set of stochastic matrices is a compact convex set with the simple and important property that  $\mathbf{A}$  is stochastic if and only if  $\mathbf{A}\mathbb{1} = \mathbb{1}$  where  $\mathbb{1}$  is the vector with all components equal to +1.

Permutation matrices are doubly stochastic matrices. If we denote by  $\mathcal{D}_n$  the set of doubly stochastic matrices, it can be proved that  $\mathcal{P}_n = O_n \cap \mathcal{D}_n$  [48]. The permutation matrices are the fundamental and prototypical doubly stochastic matrices, for Birkhoff's theorem states that any doubly stochastic matrix is a linear convex combination of finitely many permutation matrices [42]:

### Theorem 4

(Birkhoff) A matrix  $\mathbf{A}$  is a doubly stochastic matrix if and only if for some  $N < \infty$  there are permutation matrices  $\mathbf{P}_1, \dots, \mathbf{P}_N$  and positive scalars  $s_1, \dots, s_N$  such that  $s_1 + \dots + s_N = 1$  and  $\mathbf{A} = s_1\mathbf{P}_1 + \dots + s_N\mathbf{P}_N$ .

A complete proof of this theorem is to be found in [42][pages 526–528]. The proof relies on the fact that  $\mathcal{D}_n$  is a compact convex set and every point in such a set is a convex combination of the extreme points



of the set. First it is proved that every permutation matrix is an extreme point of  $\mathcal{D}_n$  and second it is shown that a given matrix is an extreme point of  $\mathcal{D}_n$  if and only if it is a permutation matrix.

## A.2 The Frobenius Norm

The Frobenius (or Euclidean) norm of a matrix  $\mathbf{A}_{n \times n}$  is an *entry-wise* norm that treats the matrix as a vector of size  $1 \times nn$ . The standard norm properties hold:  $\|\mathbf{A}\|_F > 0 \Leftrightarrow \mathbf{A} \neq 0$ ,  $\|\mathbf{A}\|_F = 0 \Leftrightarrow \mathbf{A} = 0$ ,  $\|c\mathbf{A}\|_F = c\|\mathbf{A}\|_F$ , and  $\|\mathbf{A} + \mathbf{B}\|_F \leq \|\mathbf{A}\|_F + \|\mathbf{B}\|_F$ . Additionally, the Frobenius norm is *sub-multiplicative*:

$$\|\mathbf{AB}\|_F \leq \|\mathbf{A}\|_F \|\mathbf{B}\|_F \quad (\text{A.1})$$

as well as *unitarily-invariant*. This means that for any two orthogonal matrices  $\mathbf{U}$  and  $\mathbf{V}$ :

$$\|\mathbf{UAV}\|_F = \|\mathbf{A}\|_F. \quad (\text{A.2})$$

It immediately follows the following equalities:

$$\|\mathbf{UAU}^\top\|_F = \|\mathbf{UA}\|_F = \|\mathbf{AU}\|_F = \|\mathbf{A}\|_F. \quad (\text{A.3})$$

## A.3 Spectral Properties of the Normalized Laplacian

**The normalized Laplacian** Let  $\tilde{\mathbf{u}}_k$  and  $\gamma_k$  denote the eigenvectors and eigenvalues of  $\tilde{\mathbf{L}}$ ; The spectral decomposition is  $\tilde{\mathbf{L}} = \tilde{\mathbf{U}}\mathbf{\Gamma}\tilde{\mathbf{U}}^\top$  with  $\tilde{\mathbf{U}}\tilde{\mathbf{U}}^\top = \mathbf{I}$ . The smallest eigenvalue and associated eigenvector are  $\gamma_1 = 0$  and  $\tilde{\mathbf{u}}_1 = \mathbf{D}^{1/2}\mathbb{1}$ .

We obtain the following equivalent relations:

$$\sum_{i=1}^n d_i^{1/2} \tilde{u}_{ik} = 0, \quad 2 \leq k \leq n \quad (\text{A.4})$$

$$d_i^{1/2} |\tilde{u}_{ik}| < 1, \quad 1 \leq i \leq n, 2 \leq k \leq n. \quad (\text{A.5})$$

Using (1.5) we obtain a useful expression for the combinatorial Laplacian in terms of the spectral decomposition of the normalized Laplacian. Notice, however, that the expression below is NOT a spectral decomposition of the combinatorial Laplacian:

$$\mathbf{L} = (\mathbf{D}^{1/2}\tilde{\mathbf{U}}\mathbf{\Gamma}^{1/2})(\mathbf{D}^{1/2}\tilde{\mathbf{U}}\mathbf{\Gamma}^{1/2})^\top. \quad (\text{A.6})$$

For a connected graph  $\gamma_1$  has multiplicity 1:  $0 = \gamma_1 < \gamma_2 \leq \dots \leq \gamma_n$ . As in the case of the combinatorial Laplacian, there is an upper bound on the eigenvalues (see [37] for a proof):

**Proposition 4**

For all  $k \leq n$ , we have  $\mu_k \leq 2$ .

We obtain the following spectral decomposition for the normalized Laplacian :

$$\tilde{\mathbf{L}} = \sum_{k=2}^n \gamma_k \tilde{\mathbf{u}}_k \tilde{\mathbf{u}}_k^\top. \quad (\text{A.7})$$

The spread of the graph along the  $k$ -th normalized Laplacian eigenvector is given by  $\forall(k, i), 2 \leq k \leq n, 1 \leq i \leq n$ :

$$\bar{u}_k = \frac{1}{n} \sum_{i=1}^n \tilde{u}_{ik} \quad (\text{A.8})$$

$$\sigma_{u_k} = \frac{1}{n} - \bar{u}_k^2. \quad (\text{A.9})$$

Therefore, the projection of the graph onto an eigenvector  $\tilde{\mathbf{u}}_k$  is not centered. By combining (1.5) and (A.7) we obtain an alternative representation of the combinatorial Laplacian in terms of the the spectrum of the normalized Laplacian, namely:

$$\mathbf{L} = \sum_{k=2}^n \gamma_k (\mathbf{D}^{1/2} \tilde{\mathbf{u}}_k) (\mathbf{D}^{1/2} \tilde{\mathbf{u}}_k)^\top. \quad (\text{A.10})$$

Hence, an alternative is to project the graph onto the vectors  $\mathbf{t}_k = \mathbf{D}^{1/2} \tilde{\mathbf{u}}_k$ . From  $\tilde{\mathbf{u}}_{k \geq 2}^\top \tilde{\mathbf{u}}_1 = 0$  we get that  $\mathbf{t}_{k \geq 2}^\top \mathbb{1} = 0$ . Therefore, the spread of the graph's projection onto  $\mathbf{t}_k$  has the following mean and variance,  $\forall(k, i), 2 \leq k \leq n, 1 \leq i \leq n$ :

$$\bar{t}_k = \sum_{i=1}^n d_i^{1/2} \tilde{u}_{ik} = 0 \quad (\text{A.11})$$

$$\sigma_{t_k} = \frac{1}{n} \sum_{i=1}^n d_i \tilde{u}_{ik}^2. \quad (\text{A.12})$$

**The random-walk Laplacian.** This operator is not symmetric, however its spectral properties can be easily derived from those of the normalized Laplacian using (1.7). Notice that this can be used to transform a non-symmetric Laplacian into a symmetric one, as proposed in [49] and in [50].



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